

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090822\
 Data File : BM036511.D
 Acq On : 07 Sep 2022 16:25
 Operator : CG/JU
 Sample : PB147480BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK480

Quant Time: Sep 08 00:51:30 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM081622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Sep 02 01:33:23 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.722	152	205057	20.000	ng/u1	0.00
20) Naphthalene-d8	10.516	136	918392	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.368	164	541290	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.115	188	1105238	20.000	ng/u1	0.00
79) Chrysene-d12	21.303	240	1057886	20.000	ng/u1	0.00
88) Perylene-d12	23.591	264	968871	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.128	96	29460	5.475	ng/uL	0.00
4) Pyridine-d5	3.551	84	419489	28.272	ng/u1	0.00
7) Phenol-d5	6.898	99	481419	27.505	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.063	67	347091	31.504	ng/u1	0.00
11) 2-Chlorophenol-d4	7.251	132	407741	29.649	ng/u1	0.00
15) 4-Methylphenol-d8	8.445	113	409061	31.300	ng/u1	0.00
21) Nitrobenzene-d5	8.892	128	203995	29.297	ng/u1	0.00
24) 2-Nitrophenol-d4	9.610	143	202471	28.667	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.145	165	348657	27.627	ng/u1	0.00
31) 4-Chloroaniline-d4	10.669	131	583527	29.826	ng/u1	0.00
46) Dimethylphthalate-d6	13.798	166	1126093	32.884	ng/u1	0.00
49) Acenaphthylene-d8	14.062	160	1328601	31.107	ng/u1	0.00
54) 4-Nitrophenol-d4	14.598	143	176952	27.271	ng/u1	0.00
60) Fluorene-d10	15.362	176	1043357	33.717	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.504	200	125852	22.149	ng/u1	0.00
73) Anthracene-d10	17.215	188	1648879	33.552	ng/u1	0.00
81) Pyrene-d10	19.515	212	1877385	34.211	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.450	264	1542882	32.479	ng/u1	0.00

Target Compounds	Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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